

# Can Quantum Computer Perform Better than Classical ?

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We present a theoretical model of a quantum device which can factorize any number  $N$  in two steps, by preparing an input state and then performing a proper measurement process. However, the analysis reveals that the duration of state preparation and measurement is at least proportional to  $N$  and hence the computation is not efficient. On the other hand, the energy consumption of this quantum computer grows like  $\log N$  while for classical ones is exponential in the input bit size  $m = \log_2 N$ . These results suggest the existence of a generalized Heisenberg relation which put limits on the efficiency of quantum computers in terms of the total computation time, the total energy consumption and the classical complexity of the problem.

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The action of a quantum computer is described on the abstract level in the following way. We have a quantum system with  $N$  orthogonal states (computational basis) which can store  $n = \log_2 N$  bits of information. Firstly, we have to prepare our quantum system in an initial (input) state. Then a quantum algorithm is performed which is realized as a sequence of  $T$  unitary transformation called quantum gates. Finally, the output state is measured. The number  $T$  of involved quantum gates is assumed to be proportional to the physical time needed to achieve a given task. The algorithm is called efficient if  $T$  is polynomial in  $m = \log_2 N$ . The celebrated Shor's quantum algorithm [1] which factorizes numbers into primes is efficient while its classical counterparts are not. This remarkable result had an enormous impact on the development of the whole field of quantum information and quantum computing [2].

There are, however, the following factors which can spoil the performance of a quantum computer:

A) The unavoidable inaccuracies in manufacturing of a hardware which produce deviations of the real computer's Hamiltonian from the designed one.

B) The number of elementary transformations which form a quantum algorithm is by no means unique and does not determine the real physical

time of computation. In particular, by increasing the energy level spacing of the corresponding computer's Hamiltonian we can speed up its time evolution. On the other hand the decoherence effects due to an interaction with an environment typically grow in a nonlinear way with the energy level spacing.

C) A finite duration of the input state preparation procedure and the output state measurement process should

be taken into account. In particular, one can expect that the Heisenberg energy-time uncertainty relation may put here some universal limits.

The factor A) poses a rather technological problem and will be not discussed here. The questions raised in B) will be discussed in details in the forthcoming publication [3]. Our goal is to show by constructing an explicit but purely theoretical model that there exist fundamental obstacles related to C) which could destroy the efficiency of quantum algorithms.

We describe the operation of a fictitious quantum machine which can factorize any number  $N$  with  $T = 0$  quantum gates, it means by preparing an input state and a proper measurement of it only. Although formally, there are no quantum gates in this model the quantum algorithm exists in the form of a specially designed Hamiltonian of the system. The dynamics governed by the system's Hamiltonian influences state preparation and measurements processes. In this case the problems raised in B) are neglected.

Consider a resonant cavity which supports the states of a photon (radiation modes) with the frequencies being the logarithms

of prime numbers times a fixed frequency unit  $\omega$

$$\omega_q = \omega \log q, \quad q = 2, 3, 5, 7, 11, 13, \dots \quad (1)$$

The second quantization Hamiltonian of the electromagnetic field in the cavity written in terms of annihilation and creation operators  $\hat{a}_q, \hat{a}_q^+$

$$\hat{H} = \hbar\omega \sum_q (\log q) \hat{a}_q^+ \hat{a}_q \quad (2)$$

possesses nondegenerated eigenvalues being proportional to the logarithms of all natural numbers

$$\hat{H}\psi_N = E_N\psi_N, \quad E_N = \hbar\omega \log N, \quad N = 1, 2, 3, 4, \dots \quad (3)$$

The structure of the corresponding eigenstates reveals the factorization

of  $N$  into prime numbers. Namely,

$$\psi_N \sim (\hat{a}_{q_1}^+)^{m_1} (\hat{a}_{q_2}^+)^{m_2} \dots (\hat{a}_{q_r}^+)^{m_r} \psi_1 \quad (4)$$

where

$$N = (q_1)^{m_1} (q_2)^{m_2} \dots (q_r)^{m_r} \quad (5)$$

and  $\psi_1$  is a vacuum state. In words, at the state  $\psi_N$  we have  $m_1$  photons of the frequency  $\omega \log q_1$ ,  $m_2$  photons of the frequency  $\omega \log q_2$ , ..., and  $m_r$  photons of the

frequency  $\omega \log q_r$ . Therefore, in principle one can find the factorization of any number  $N$  in two steps. First we prepare the system in the state  $\psi_N$  of a given energy  $E_N = \hbar\omega \log N$  by transferring this portion of energy into the empty cavity. Then we open the cavity and perform a spectral analysis of the corresponding radiation field counting photons in different modes.

Let us discuss the possible preparation process. We perturb our quantum system being initially in a vacuum state by a weak external interaction Hamiltonian  $\hat{V}(t)$  which is periodic in time with a tunable frequency  $\Omega$

$$\hat{V}(t) = \hat{W} \cos(\Omega t) . \quad (6)$$

Instead of a usual electromagnetic interaction linear in field we assume that  $\hat{W}$  is a sufficiently nonlinear function of electromagnetic field operators which allows multiphoton excitations such that the (virtual) transitions between the vacuum  $\psi_1$  and any state  $\psi_N$  are possible i.e.

$$\langle \psi_1, \hat{W} \psi_N \rangle \neq 0 , \text{ for all } N = 2, 3, 4, \dots \quad (7)$$

For a given number  $N$  which we want to factorize we tune the frequency  $\Omega$  to the value  $\omega \log N$ . The time dependent first-order perturbation calculus [4] gives us the probability of excitation of the state  $\psi_M$

$$p_M(t) = \frac{2}{\hbar^2} |\langle \psi_1, \hat{W} \psi_M \rangle|^2 \frac{\sin^2 \left\{ \frac{1}{2} \omega (\log M - \log N) t \right\}}{\omega^2 (\log M - \log N)^2} . \quad (8)$$

As the energy level spacing around  $E_N = \hbar\omega \log N$  is  $\delta E_N \approx \hbar\omega/N$  it follows from the formula (8) that we have to wait for a time at least of the order

$$t \approx N\omega^{-1} \quad (9)$$

to be sure that the selected state  $\psi_N$  has been prepared with much larger probability than the other neighboring states  $\psi_M$ . The similar estimation can be easily obtained for the duration

of the measurement process. Therefore the total computation time  $t_c$  grows exponentially with  $\log N$  like in the classical situation.

It is obvious that the result obtained above can be treated as a special case of the Heisenberg time-energy uncertainty relation [4]

$$\Delta t \cdot \Delta E \geq \frac{\hbar}{2} . \quad (10)$$

Indeed in order to identify the energy of a quantum state with an accuracy  $\hbar\omega/N$  we need a time longer than  $N/\omega$ .

One should notice, however, that our quantum computer is superior to classical ones in respect of energy consumption, at least for the case of existing irreversible computers (see [5] for the theory of reversible computations). The energy used for the factorization of  $N$  is

equal to  $E_c = \hbar\omega \log N$  while for the classical irreversible computers any logical step consumes an energy portion and hence the total energy cost of factorization grows exponentially with the input bit size  $m = \log_2 N$ . Taking into account the eq.(9) we obtain the following inequality independent of an arbitrary frequency scale  $\omega$

$$t_c E_c \gg \hbar N \log N . \quad (11)$$

The form of the inequality (11) suggests the following general hypothesis. There exists an inequality which puts universal limits on the performance of a quantum computer in terms of the total computation time  $t_c$ , the total energy consumption  $E_c$  and the complexity  $\mathcal{C}(\log_2 N)$  of the problem to be solved. This "generalized Heisenberg relation" reads

$$t_c E_c \gg \hbar \mathcal{C}(\log_2 N) . \quad (12)$$

The complexity  $\mathcal{C}(\log_2 N)$  is a function of the input bit size and is defined by a minimal number of logical steps needed to solve the problem. The inequality (12) means that for a non-efficient optimal classical algorithm the quantum computation is also not efficient either with respect to the computation time or the used energy.

In order to prove this hypothesis we cannot restrict ourselves to counting quantum gates in the algorithm but we have to discuss physical implementation of all steps of quantum computing including state preparation and measurement processes.

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